Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112116\
Data File : BF091265.D
Acq On : 22 Nov 2016 1:53
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 22 10:59:04 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 22 10:25:16 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	99	0.00
2		1,4-Dioxane	0.547	0.529	3.3	98	0.00
3		Pyridine	1.430	1.605	-12.2	101	-0.01
4		n-Nitrosodimethylamine	0.772	0.746	3.4	91	0.00
5	S	2-Fluorophenol	1.203	1.198	0.4	99	0.00
6		Aniline	1.728	1.795	-3.9	100	0.00
7	S	Phenol-d6	1.439	1.496	-4.0	100	0.00
8		2-Chlorophenol	1.289	1.261	2.2	99	-0.01
9		Benzaldehyde	0.722	0.783	-8.4	104	0.00
10	C	Phenol	1.683	1.662	1.2	95	0.00
11		bis(2-Chloroethyl)ether	1.311	1.300	0.8	99	0.00
12		1,3-Dichlorobenzene	1.401	1.391	0.7	97	0.00
13	C	1,4-Dichlorobenzene	1.464	1.448	1.1	98	0.00
14		1,2-Dichlorobenzene	1.300	1.285	1.2	98	0.00
15		Benzyl Alcohol	1.130	1.160	-2.7	98	0.00
16		2,2'-oxybis(1-Chloropropane	2.510	2.265	9.8	92	0.00
17		2-Methylphenol	1.296	1.378	-6.3	99	0.00
18		Hexachloroethane	0.542	0.534	1.5	97	0.00
19	P	n-Nitroso-di-n-propylamine	1.051	0.976	7.1	99	0.00
20		3+4-Methylphenols	1.296	1.378	-6.3	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
22		Acetophenone	0.477	0.476	0.2	100	0.00
23	S	Nitrobenzene-d5	0.363	0.364	-0.3	96	0.00
24		Nitrobenzene	0.404	0.386	4.5	99	0.00
25		Isophorone	0.684	0.663	3.1	96	0.00
26	C	2-Nitrophenol	0.172	0.178	-3.5	96	0.00
27		2,4-Dimethylphenol	0.272	0.279	-2.6	99	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.393	-5.1	99	0.00
29	C	2,4-Dichlorophenol	0.270	0.282	-4.4	96	0.00
30		1,2,4-Trichlorobenzene	0.302	0.289	4.3	96	0.00
31		Naphthalene	0.934	0.939	-0.5	98	0.00
32		Benzoic acid	0.158	0.184	-16.5	94	0.00
33		4-Chloroaniline	0.378	0.387	-2.4	99	0.00
34	C	Hexachlorobutadiene	0.179	0.174	2.8	98	0.00
35		Caprolactam	0.076	0.075	1.3	100	0.00
36	C	4-Chloro-3-methylphenol	0.296	0.300	-1.4	99	0.00
37		2-Methylnaphthalene	0.627	0.610	2.7	98	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
39		1,2,4,5-Tetrachlorobenzene	0.564	0.530	6.0	100	0.00
40	P	Hexachlorocyclopentadiene	0.125	0.139	-11.2	90	0.00
41	S	2,4,6-Tribromophenol	0.187	0.176	5.9	101	0.00
42	C	2,4,6-Trichlorophenol	0.343	0.329	4.1	96	0.00
43		2,4,5-Trichlorophenol	0.360	0.345	4.2	101	0.00
44	S	2-Fluorobiphenyl	1.145	1.196	-4.5	99	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.525	1.332	12.7	96	0.00
46		2-Chloronaphthalene	1.135	1.047	7.8	96	0.00
47		2-Nitroaniline	0.421	0.411	2.4	98	0.00
48		Acenaphthylene	1.708	1.635	4.3	98	0.00
49		Dimethylphthalate	1.334	1.316	1.3	99	0.00
50		2,6-Dinitrotoluene	0.286	0.287	-0.3	101	0.00
51	C	Acenaphthene	1.059	1.035	2.3	100	0.00
52		3-Nitroaniline	0.305	0.337	-10.5	102	0.00
53	P	2,4-Dinitrophenol	0.146	0.132	9.6	88	0.00
54		Dibenzofuran	1.475	1.374	6.8	97	0.00
55	P	4-Nitrophenol	0.123	0.132	-7.3	98	0.00
56		2,4-Dinitrotoluene	0.360	0.330	8.3	99	-0.01
57		Fluorene	1.194	1.178	1.3	100	0.00
58		2,3,4,6-Tetrachlorophenol	0.303	0.297	2.0	99	0.00
59		Diethylphthalate	1.294	1.137	12.1	97	0.00
60		4-Chlorophenyl-phenylether	0.561	0.523	6.8	100	0.00
61		4-Nitroaniline	0.314	0.333	-6.1	102	0.00
62		Azobenzene	1.538	1.453	5.5	102	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
64		4,6-Dinitro-2-methylphenol	0.128	0.115	10.2	93	0.00
65	c	n-Nitrosodiphenylamine	0.603	0.569	5.6	100	0.00
66		4-Bromophenyl-phenylether	0.209	0.206	1.4	103	0.00
67		Hexachlorobenzene	0.228	0.229	-0.4	101	0.00
68		Atrazine	0.168	0.163	3.0	102	0.00
69	C	Pentachlorophenol	0.090	0.089	1.1	96	0.00
70		Phenanthrene	1.051	0.931	11.4	99	0.00
71		Anthracene	1.047	1.101	-5.2	107	0.00
72		Carbazole	0.960	0.899	6.4	100	0.00
73		Di-n-butylphthalate	1.142	1.178	-3.2	102	0.00
74	C	Fluoranthene	1.145	1.109	3.1	99	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76		Benzidine	0.449	0.384	14.5	78	0.00
77		Pyrene	1.534	1.395	9.1	96	-0.01
78	S	Terphenyl-d14	0.882	0.878	0.5	102	0.00
79		Butylbenzylphthalate	0.644	0.688	-6.8	103	0.00
80		Benzo(a)anthracene	1.212	1.160	4.3	102	0.00
81		3,3'-Dichlorobenzidine	0.420	0.432	-2.9	99	0.00
82		Chrysene	1.246	1.309	-5.1	106	0.00
83		Bis(2-ethylhexyl)phthalate	0.816	0.814	0.2	103	0.00
84	c	Di-n-octyl phthalate	1.429	1.441	-0.8	105	-0.01
85		Indeno(1,2,3-cd)pyrene	0.991	0.977	1.4	107	0.00
86	I	Perylene-d12	1.000	1.000	0.0	103	0.00
87		Benzo(b)fluoranthene	1.202	1.144	4.8	106	0.00

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88	Benzo(k)fluoranthene	1.214	1.081	11.0	101	0.00
89 C	Benzo(a)pyrene	1.102	1.064	3.4	103	0.00
90	Dibenzo(a,h)anthracene	0.908	0.948	-4.4	107	0.00
91	Benzo(g,h,i)perylene	0.928	0.974	-5.0	108	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0